```
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PASSWORD:
TERMINAL (ENTER 1, 2, 3, OR ?):2


* * * * * * * * * *   Welcome to STN International   * * * * * * * * * *

 NEWS  1           Web Page URLs for STN Seminar Schedule - N. America
 NEWS  2           "Ask CAS" for self-help around the clock
 NEWS  3  JUL 12   BEILSTEIN enhanced with new display and select options,
                   resulting in a closer connection to BABS
 NEWS  4  AUG 02   IFIPAT/IFIUDB/IFICDB reloaded with new search and display
                   fields
 NEWS  5  AUG 02   CAplus and CA patent records enhanced with European and Japan
                   Patent Office Classifications
 NEWS  6  AUG 02   The Analysis Edition of STN Express with Discover!
                   (Version 7.01 for Windows) now available
 NEWS  7  AUG 27   BIOCOMMERCE: Changes and enhancements to content coverage
 NEWS  8  AUG 27   BIOTECHABS/BIOTECHDS: Two new display fields added for legal
                   status data from INPADOC
 NEWS  9  SEP 01   INPADOC: New family current-awareness alert (SDI) available
 NEWS 10  SEP 01   New pricing for the Save Answers for SciFinder Wizard within
                   STN Express with Discover!
 NEWS 11  SEP 01   New display format, HITSTR, available in WPIDS/WPINDEX/WPIX
 NEWS 12  SEP 27   STANDARDS will no longer be available on STN
 NEWS 13  SEP 27   SWETSCAN will no longer be available on STN
 NEWS 14  OCT 28   KOREAPAT now available on STN

 NEWS EXPRESS  OCTOBER 29 CURRENT WINDOWS VERSION IS V7.01A, CURRENT
               MACINTOSH VERSION IS V6.0c(ENG) AND V6.0Jc(JP),
               AND CURRENT DISCOVER FILE IS DATED 11 AUGUST 2004
 NEWS HOURS    STN Operating Hours Plus Help Desk Availability
 NEWS INTER    General Internet Information
 NEWS LOGIN    Welcome Banner and News Items
 NEWS PHONE    Direct Dial and Telecommunication Network Access to STN
 NEWS WWW      CAS World Wide Web Site (general information)

Enter NEWS followed by the item number or name to see news on that
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  of commercial gateways or other similar uses is prohibited and may
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  * * * * * * * * * * * * * * STN  Columbus  * * * * * * * * * * * * * * *

FILE 'HOME' ENTERED AT 14:38:48 ON 09 NOV 2004

=> fil reg
FILE 'REGISTRY' ENTERED AT 14:39:05 ON 09 NOV 2004
USE IS SUBJECT TO THE TERMS OF YOUR STN CUSTOMER AGREEMENT.
PLEASE SEE "HELP USAGETERMS" FOR DETAILS.
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```

Property values tagged with IC are from the ZIC/VINITI data file
provided by InfoChem.

STRUCTURE FILE UPDATES: 8 NOV 2004 HIGHEST RN 777024-10-9
DICTIONARY FILE UPDATES: 8 NOV 2004 HIGHEST RN 777024-10-9

TSCA INFORMATION NOW CURRENT THROUGH MAY 21, 2004

Please note that search-term pricing does apply when
conducting SmartSELECT searches.

Crossover limits have been increased. See HELP CROSSOVER for details.

Experimental and calculated property data are now available. For more
information enter HELP PROP at an arrow prompt in the file or refer
to the file summary sheet on the web at:
http://www.cas.org/ONLINE/DBSS/registryss.html

*** YOU HAVE NEW MAIL ***
'.REGISTRY' IS DEFAULT FORMAT FOR 'REGISTRY' FILE

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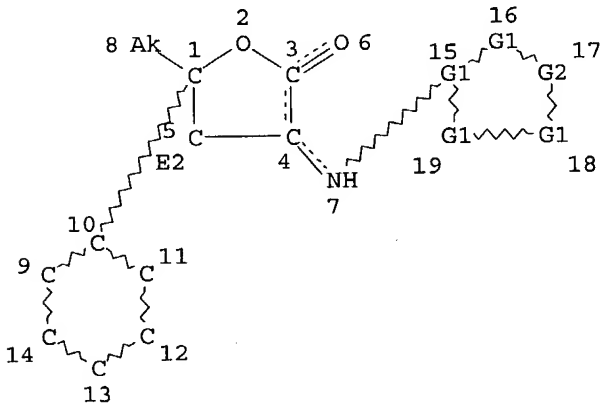
L1 STR



VAR G1=C/N
REP G2=(1-2) C
NODE ATTRIBUTES:
HCOUNT IS E2 AT 5
DEFAULT MLEVEL IS ATOM
DEFAULT ECLEVEL IS LIMITED

GRAPH ATTRIBUTES:
RING(S) ARE ISOLATED OR EMBEDDED
NUMBER OF NODES IS 19

STEREO ATTRIBUTES: NONE

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SAMPLE SEARCH INITIATED 14:56:43 FILE 'REGISTRY'
SAMPLE SCREEN SEARCH COMPLETED - 67 TO ITERATE

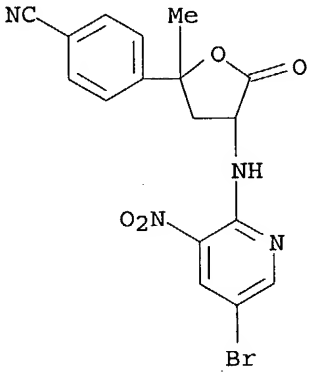
100.0% PROCESSED 67 ITERATIONS 6 ANSWERS
SEARCH TIME: 00.00.01

FULL FILE PROJECTIONS: ONLINE **COMPLETE**
BATCH **COMPLETE**
PROJECTED ITERATIONS: 849 TO 1831
PROJECTED ANSWERS: 6 TO 266

L2 6 SEA SSS SAM L1

=> d scan

L2 6 ANSWERS REGISTRY COPYRIGHT 2004 ACS on STN
IN Pentonic acid, 2-[(5-bromo-3-nitro-2-pyridinyl)amino]-4-C-(4-cyanophenyl)-2,3,5-trideoxy-, γ-lactone (9CI)
MF C17 H13 Br N4 O4



PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

HOW MANY MORE ANSWERS DO YOU WISH TO SCAN? (1):0

=> s l1 ful
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FULL SCREEN SEARCH COMPLETED - 1383 TO ITERATE

100.0% PROCESSED 1383 ITERATIONS 172 ANSWERS
SEARCH TIME: 00.00.01

L3 172 SEA SSS FUL L1

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172	RN	325482-22-2	REGISTRY

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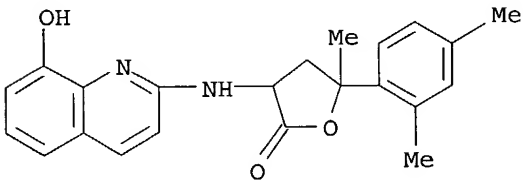
L3 ANSWER 68 OF 172 REGISTRY COPYRIGHT 2004 ACS on STN
RN 486447-75-0 REGISTRY
CN Pentonic acid, 2,3,5-trideoxy-4-C-(2,4-dimethylphenyl)-2-[(8-hydroxy-2-quinolinyl)amino]-, γ-lactone (9CI) (CA INDEX NAME)
FS 3D CONCORD
MF C22 H22 N2 O3
SR CA
LC STN Files: CA, CAPLUS, TOXCENTER, USPATFULL
DT.CA CAplus document type: Patent
RL.P Roles from patents: BIOL (Biological study); PREP (Preparation); USES (Uses)



PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

1 REFERENCES IN FILE CA (1907 TO DATE)
1 REFERENCES IN FILE CAPLUS (1907 TO DATE)

REFERENCE 1

AN 138:106595 CA
TI Use of substituted gamma-lactone compounds as medicaments
IN Maul, Corinna; Sundermann, Bernd; Przewosny, Michael; Hennies, Hagen-Heinrich
PA Gruenenthal G.m.b.H., Germany
SO PCT Int. Appl., 63 pp.
CODEN: PIXXD2
DT Patent
LA German
FAN.CNT 1

PATENT NO.	KIND	DATE	APPLICATION NO.	DATE

PI WO 2003004016 A1 20030116 WO 2002-EP7382 20020703
WO 2003004016 C1 20030213
W: AE, AG, AL, AM, AT, AU, AZ, BA, BB, BG, BR, BY, BZ, CA, CH, CN, CO, CR, CU, CZ, DK, DM, DZ, EC, EE, ES, FI, GB, GD, GE, GH, GM, HR, HU, ID, IL, IN, IS, JP, KE, KG, KP, KR, KZ, LC, LK, LR, LS, LT, LU, LV, MA, MD, MG, MK, MN, MW, MX, MZ, NO, NZ, OM, PH, PL, PT, RO, RU, SD, SE, SG, SI, SK, SL, TJ, TM, TN, TR, TT, TZ, UA, UG, US, UZ, VN, YU, ZA, ZM, ZW, AM, AZ, BY, KG, KZ, MD, RU, TJ, TM
RW: GH, GM, KE, LS, MW, MZ, SD, SL, SZ, TZ, UG, ZM, ZW, AT, BE, BG, CH, CY, CZ, DE, DK, EE, ES, FI, FR, GB, GR, IE, IT, LU, MC, NL, PT, SE, SK, TR, BF, BJ, CF, CG, CI, CM, GA, GN, GQ, GW, ML, MR, NE, SN, TD, TG
DE 10132726 A1 20030227 DE 2001-10132726 20010705
EP 1406610 A1 20040414 EP 2002-754819 20020703
R: AT, BE, CH, DE, DK, ES, FR, GB, GR, IT, LI, LU, NL, SE, MC, PT, IE, SI, LT, LV, FI, RO, MK, CY, AL, TR, BG, CZ, EE, SK
US 2004176394 A1 20040909 US 2004-751737 20040105
PRAI DE 2001-10132726 20010705
WO 2002-EP7382 20020703
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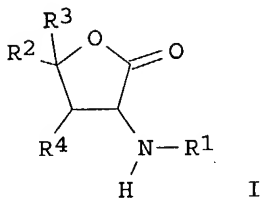
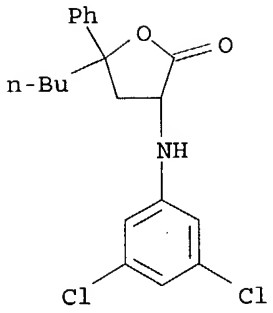


I

AB Substituted gamma-lactone compds. [I; R1 = (un)substituted 2-pyridyl, (un)substituted 2-pyrimidyl-, (un)substituted 2-quinolinyl, (un)substituted 2-pyrazinyl; R2 = (un)substituted (un)branched (un)saturated C1-10 aliphatic residue; R3 = (un)substituted aryl; R4 = H; etc.; e.g., 3-(4,6-dimethylpyridin-2-ylamino)-5-methyl-5-p-tolyldihydrofuran-2-one which demonstrated a 60% reduction of nitrogen monoxide synthase activity at 10 μM], which are prepared in an automated device, are useful in the production of medicines for treating septic shock, neurodegenerative diseases, multiple sclerosis, Parkinsonism, Alzheimer's disease, Huntington's disease, inflammation and related pain, cerebral ischemia, diabetes, meningitis, arteriosclerosis, cancer, mycosis, or for promoting wound healing.
RE.CNT 4 THERE ARE 4 CITED REFERENCES AVAILABLE FOR THIS RECORD
ALL CITATIONS AVAILABLE IN THE RE FORMAT

L3 ANSWER 69 OF 172 REGISTRY COPYRIGHT 2004 ACS on STN
RN 484053-57-8 REGISTRY
CN 2(3H)-Furanone, 5-butyl-3-[(3,5-dichlorophenyl)amino]dihydro-5-phenyl- (9CI) (CA INDEX NAME)
MF C20 H21 Cl2 N O2
SR CA
LC STN Files: CA, CAPLUS, USPATFULL
DT.CA CAplus document type: Patent
RL.P Roles from patents: BIOL (Biological study); PREP (Preparation); USES (Uses)



PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

1 REFERENCES IN FILE CA (1907 TO DATE)
1 REFERENCES IN FILE CAPLUS (1907 TO DATE)

REFERENCE 1

AN 138:89684 CA
TI Preparation of γ-lactones as NMDA antagonists
IN Maul, Corinna; Przewosny, Michael; Englberger, Werner
PA Gruenenthal G.m.b.H., Germany
SO PCT Int. Appl., 66 pp.
CODEN: PIXXD2
DT Patent
LA German
FAN.CNT 1

	PATENT NO.	KIND	DATE	APPLICATION NO.	DATE
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PRAI	DE 2001-10132725		20010705		
	WO 2002-EP7380		20020703		

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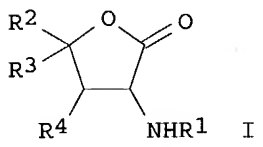


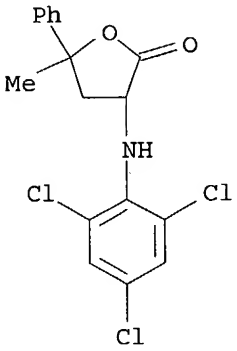
I

AB Title compds. [I; R1 = (substituted) (hetero)aryl, (hetero)arylalkyl, (substituted) (saturated) (branched) (cyclo)aliphatic group; R2 = (substituted) (saturated) (branched) aliphatic group; R3 = (substituted) aryl; R4 = H; or R3R4

= (substituted) (saturated) aliphatic residue if R2 = (substituted) aryl, (substituted) (saturated) aliphatic residue], were prepared Several I at 10 μM showed affinity to glycine binding site of NMDA receptor channel with IC50 = 62-80% (Ki = 0.8-1.3 μM).
RE.CNT 6 THERE ARE 6 CITED REFERENCES AVAILABLE FOR THIS RECORD
ALL CITATIONS AVAILABLE IN THE RE FORMAT

=> d 172 sub bib abs

L3 ANSWER 172 OF 172 REGISTRY COPYRIGHT 2004 ACS on STN
RN 325482-22-2 REGISTRY
CN Pentonic acid, 2,3,5-trideoxy-4-C-phenyl-2-[(2,4,6-trichlorophenyl)amino]-, γ-lactone (9CI) (CA INDEX NAME)
FS 3D CONCORD
MF C17 H14 Cl3 N O2
SR CA
LC STN Files: CA, CAPLUS, CASREACT
DT.CA CAplus document type: Journal
RL.NP Roles from non-patents: PREP (Preparation)



PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

1 REFERENCES IN FILE CA (1907 TO DATE)
1 REFERENCES IN FILE CAPLUS (1907 TO DATE)

REFERENCE 1

AN 134:162877 CA
TI Synthesis of α-amino γ-lactone via a novel tandem three-component reaction of alkenes, glyoxylates and amines
AU Huang, T.; Li, C.-J.
CS Department of Chemistry, Tulane University, New Orleans, LA, 70118, USA
SO Tetrahedron Letters (2000), 41(50), 9747-9751
CODEN: TELEAY; ISSN: 0040-4039
PB Elsevier Science Ltd.
DT Journal
LA English
AB α-Amino γ-lactones were generated by an InCl3-mediated or Sc(O3SCF3)3-catalyzed three-component reaction of alkenes, glyoxylates, and amines.
RE.CNT 33 THERE ARE 33 CITED REFERENCES AVAILABLE FOR THIS RECORD
ALL CITATIONS AVAILABLE IN THE RE FORMAT

=> fil caplus
FILE 'CAPLUS' ENTERED AT 14:59:52 ON 09 NOV 2004

FILE COVERS 1907 - 9 Nov 2004 VOL 141 ISS 20
FILE LAST UPDATED: 8 Nov 2004 (20041108/ED)

This file contains CAS Registry Numbers for easy and accurate substance identification.

=> s l3
L4 3 L3

=> fil beil
FILE 'BEILSTEIN' ENTERED AT 15:00:30 ON 09 NOV 2004

FILE RELOADED ON OCTOBER 20, 2002
FILE LAST UPDATED ON NOVEMBER 3, 2004

FILE COVERS 1771 TO 2004.
*** FILE CONTAINS 9,073,068 SUBSTANCES ***

>>>PLEASE NOTE: Reaction Data and substance data are stored in separate documents and can not be searched together in one query. Reaction data for BEILSTEIN compounds may be displayed immediately with the display codes PRE (preparations) and REA (reactions). A substance answer set retrieved after the search for a chemical name, a compounds with available reaction information by combining with PRE/FA, REA/FA or more generally with RX/FA. The BEILSTEIN Registry Number (BRN) is the link between a BEILSTEIN compound and belonging reactions. For mo detailed reaction searches BRNs can be searched as reaction partner BRNs Reactant BRN (RX.RBRN) or Product BRN (RX.PBRN).<<<

>>> FOR SEARCHING PREPARATIONS SEE HELP PRE <<<

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* PLEASE NOTE THAT THERE ARE NO FORMATS FREE OF COST. *
* SET NOTICE FEATURE: THE COST ESTIMATES CALCULATED FOR SET NOTICE *
* ARE BASED ON THE HIGHEST PRICE CATEGORY. THEREFORE; THESE *
* ESTIMATES MAY NOT REFLECT THE ACTUAL COSTS. *
* FOR PRICE INFORMATION SEE HELP COST *
****

NEW

*** PATENT NUMBERS (PN) AND BABS ACCESSION NUMBERS (BABSAN) CAN NOW BE SEARCHED, SELECTED AND TRANSFERRED.**
*** NEW DISPLAY FORMATS ALLREF, ALLP AND BABSAN SHOW ALL REFERENCES, ALL PATENT REFERENCES, OR ALL BABS ACCESSION NUMBERS FOR A COMPOUND AT A GLANCE.**

*** YOU HAVE NEW MAIL ***

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=> s l1 ful
FULL SEARCH INITIATED 15:00:36 FILE 'BEILSTEIN'
FULL SCREEN SEARCH COMPLETED -    674 TO ITERATE

100.0% PROCESSED     674 ITERATIONS                          1 ANSWERS
SEARCH TIME: 00.00.10

L5            1 SEA SSS FUL L1

=> d ide
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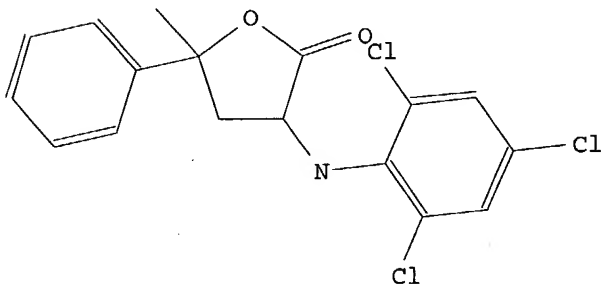
L5 ANSWER 1 OF 1 BEILSTEIN COPYRIGHT 2004 BEILSTEIN MDL on STN

Beilstein Records (BRN):	8708878
Chemical Name (CN):	5-methyl-5-phenyl-3-(2,4,6-trichloro-phenylamino)-dihydro-furan-2-one
Autonom Name (AUN):	5-methyl-5-phenyl-3-(2,4,6-trichloro-phenylamino)-dihydro-furan-2-one
Molec. Formula (MF):	C17 H14 Cl3 N O2
Molecular Weight (MW):	370.66
Lawson Number (LN):	20596, 14134
Compound Type (CTYPE):	heterocyclic
Constitution ID (CONSID):	7376333
Tautomer ID (TAUTID):	8207212
Entry Date (DED):	2001/04/26
Update Date (DUPD):	2001/04/26



Field Availability:

Code	Name	Occurrence
BRN	Beilstein Records	1
CN	Chemical Name	1
AUN	Autonomname	1
MF	Molecular Formula	1
FW	Formular Weight	1
LN	Lawson Number	2
FS	File Segment	1
CTYPE	Compound Type	1
CONSID	Constitution ID	1
TAUTID	Tautomer ID	1
ED	Entry Date	1
UPD	Update Date	1

This substance also occurs in Reaction Documents:

Code	Name	Occurrence

RX	Reaction Documents	1
RXPRO	Substance is Reaction Product	1

=> log h

COST IN U.S. DOLLARS	SINCE FILE ENTRY	TOTAL SESSION
FULL ESTIMATED COST	9.33	222.41

DISCOUNT AMOUNTS (FOR QUALIFYING ACCOUNTS)	SINCE FILE ENTRY	TOTAL SESSION
CA SUBSCRIBER PRICE	0.00	-1.98

SESSION WILL BE HELD FOR 60 MINUTES
STN INTERNATIONAL SESSION SUSPENDED AT 15:01:28 ON 09 NOV 2004